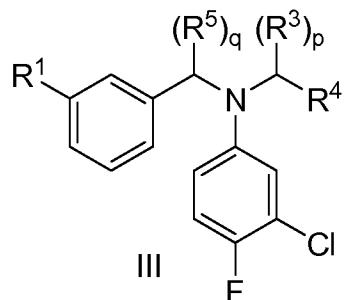


IN THE CLAIMS:

1. Cancelled
2. Cancelled
3. (Currently amended) The compound according to ~~Claim 2~~ of the Formula III;



wherein:

R¹ is selected from: H, F and OH;

R³ is selected from: H, (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, (C₂-C₆)alkenyl and (C₁-C₆)alkyl hydroxyl;

R⁴ is selected from: H, oxo, OH, halo, CN, NH₂, NO₂, (C=O)_aO_b(C₁-C₁₀)alkyl, (C=O)_aO_b(C₂-C₁₀)alkenyl, (C=O)_aO_b(C₂-C₁₀)alkynyl, (C=O)_aO_b(C₃-C₆)cycloalkyl, (C=O)_aO_b(C₀-C₆)alkylene-aryl, (C=O)_aO_b(C₀-C₆)alkylene-heterocyclyl, (C=O)_aO_b(C₀-C₆)alkylene-N(R^b)₂, C(O)H, (C₀-C₆)alkylene-CO₂H, C(O)N(R^b)₂, and S(O)₂N(R^b)₂; said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R^b;

R⁵ is H or CH₃;

R^a is selected from: (C₁-C₆)alkyl, (C₃-C₆)cycloalkyl, aryl and heterocyclyl; said alkyl, cycloalkyl, aryl and heterocyclyl is optionally substituted with one or more substituents selected from OH, (C₁-C₆)alkyl, (C₁-C₆)alkoxy, halogen, CO₂H, CN, (O)C=O(C₁-C₆)alkyl, oxo and N(R^c)₂;

R_b is independently selected from: H, oxo, OH, halogen, CO₂H, CN, (O)C=O(C₁-C₆)alkyl, N(R^c)₂, (C₁-C₆)alkyl, aryl, heterocyclyl, (C₃-C₆)cycloalkyl, (C=O)O(C₁-C₆)alkyl, C=O(C₁-C₆)alkyl and S(O)₂R^a; said alkyl, cycloalkyl, aryl or heterocyclyl is optionally substituted with one or more substituents selected from OH, (C₁-C₆)alkyl, (C₁-C₆)alkoxy, halogen, CO₂H, CN, (O)C=O(C₁-C₆)alkyl, oxo, N(R^c)₂ and optionally substituted heterocyclyl, wherein said heterocyclyl is optionally substituted with (C₁-C₆)alkyl, oxo or NH₂.

R^c is independently selected from: H and (C₁-C₆)alkyl;

a is 0 or 1;

b is 0 or 1;

p is 1 or 2;

q is 0 or 1;

~~and all other substituents and variables are as defined in Claim 2;~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

4. (Original) The compound according to Claim 3 of the Formula III;

wherein:

R⁴ is selected from: H, oxo, OH, halo, CN, NH₂, NO₂, (C=O)aOb(C₁-C₁₀)alkyl, (C=O)aOb(C₂-C₁₀)alkenyl, (C=O)aOb(C₂-C₁₀)alkynyl, (C=O)aOb(C₃-C₆)cycloalkyl, (C=O)aOb(C₀-C₆)alkylene-aryl, (C=O)aOb(C₀-C₆)alkylene-heterocyclyl, (C=O)aOb(C₀-C₆)alkylene-N(R^b)₂, C(O)H, (C₀-C₆)alkylene-CO₂H, C(O)N(R^b)₂, and S(O)₂N(R^b)₂; said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R^b;

and all other substituents and variables are as defined in Claim 3;

or a pharmaceutically acceptable salt or stereoisomer thereof.

5. (Original) The compound according to Claim 3 of the Formula III;

wherein:

R⁴ is selected from: H, oxo, OH, halo, CN, NH₂, NO₂, (C=O)_aO_b(C₁-C₁₀)alkyl, (C=O)_aO_b(C₂-C₁₀)alkenyl, (C=O)_aO_b(C₂-C₁₀)alkynyl, (C=O)_aO_b(C₀-C₆)alkylene-N(R^b)₂, (C=O)-R^b, C(O)H, (C₀-C₆)alkylene-CO₂H, C(O)N(R^b)₂, and S(O)₂N(R^b)₂; said alkyl, alkenyl, alkynyl and alkylene is optionally substituted with up to three substituents selected from R^b;

and all other substituents and variables are as defined in Claim 3;

or a pharmaceutically acceptable salt or stereoisomer thereof.

6. (Original) A compound which is selected from:

3-{[(3-chloro-4-fluorophenyl)(2-hydroxy-1-methylethyl)amino]methyl}phenol;
2-[benzyl(3-chloro-4-fluorophenyl)amino]propan-1-ol;
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1-[2(dimethylamino)ethyl] alaninamide;
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)alaninamide;
methyl N-benzyl-N-(3-chloro-4-fluorophenyl)alanylglycinate;
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1-(isoxazol-4-ylmethyl)alaninamide;
3-[benzyl(3-chloro-4-fluorophenyl)amino]-2-methylbutan-2-ol;
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1,N-1-dimethylpropane-1,2-diamine;
N-benzyl-3-chloro-4-fluoro-N-[1-methyl-2-(4-methylpiperazin-1-yl)ethyl]aniline;
2-[(3-chloro-4-fluorophenyl)(1-phenylethyl)amino]propan-1-ol;
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)-N-1-(isoxazol-4-ylmethyl)alaninamide;
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)alaninamide;
N-2-(3-chloro-4-fluorophenyl)-N-1-[2-(dimethylamino)ethyl]-N-2-(3-hydroxybenzyl)alaninamide;
Methyl 2-[benzyl(3-chloro-4-fluorophenyl)amino]butanoate;
Methyl 2-[benzyl(3-chloro-4-fluorophenyl)amino]pent-4-enoate;
2-[benzyl(3-chloro-4-fluorophenyl)amino]pent-4-en-1-ol;
N-benzyl-N-(3-chloro-4-fluorophenyl)glycine;
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentan-1-ol;
2-[benzyl(3-chloro-4-fluorophenyl)amino]butan-1-ol;
N-benzyl-3-chloro-N-[1-(3-[(dimethylamino)methyl]piperidin-1-yl)carbonyl]propyl]-4-fluoroaniline;
2-[benzyl(3-chloro-4-fluorophenyl)amino]-N-methyl-N-[2-(1-methyl-1H-pyrazol-4-yl)ethyl]butanamide;

2-[benzyl(3-chloro-4-fluorophenyl)amino]-3-methylbutan-1-ol;
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentane-1,5-diol;
2-[benzyl(3-chloro-4-fluorophenyl)amino]-3-cyclopropylpropan-1-ol; and
N²-benzyl-N²-(3-chloro-4-fluorophenyl)-N¹-[2-(dimethylamino)ethyl]-2-methylalaninamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

7. (Currently amended) The TFA salt of a compound according to Claim 43 which is

3-{[(3-chloro-4-fluorophenyl)(2-hydroxy-1-methylethyl)amino]methyl}phenol;
2-[benzyl(3-chloro-4-fluorophenyl)amino]propan-1-ol;
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)-N-1-[2(dimethylamino)ethyl] alaninamide;
N-2-benzyl-N-2-(3-chloro-4-fluorophenyl)alaninamide;
methyl N-benzyl-N-(3-chloro-4-fluorophenyl)alanyl glycinate;
3-[benzyl(3-chloro-4-fluorophenyl)amino]-2-methylbutan-2-ol;
2-[3-chloro-4-fluorophenyl](1-phenylethyl)amino]propan-1-ol;
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)-N-1-(isoxazol-4-ylmethyl)alaninamide;
N-2-(3-chloro-4-fluorophenyl)-N-2-(3-hydroxybenzyl)alaninamide;
N-2-(3-chloro-4-fluorophenyl)-N-1-[2-(dimethylamino)ethyl]-N-2-(3-hydroxybenzyl)alaninamide;
N-benzyl-N-(3-chloro-4-fluorophenyl)glycine;
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentan-1-ol;
2-[benzyl(3-chloro-4-fluorophenyl)amino]butan-1-ol;
N-benzyl-3-chloro-N-[1-{3-[(dimethylamino)methyl]piperidin-1-yl}carbonyl]propyl]-4-fluoroaniline;
2-[benzyl(3-chloro-4-fluorophenyl)amino]-N-methyl-N-[2-(1-methyl-1H-pyrazol-4-yl)ethyl]butanamide;
2-[benzyl(3-chloro-4-fluorophenyl)amino]pentane-1,5-diol; and
N²-benzyl-N²-(3-chloro-4-fluorophenyl)-N¹-[2-(dimethylamino)ethyl]-2-methylalaninamide;

or stereoisomer thereof.

8. (Currently amended) A pharmaceutical composition comprising a pharmaceutical carrier, and dispersed therein, a therapeutically effective amount of a compound of Claim 34.

9. (Currently amended) ~~The use of the compound according to Claim 1 for the preparation of a medicament useful in the treatment or prevention of~~ A method of treating cancer in a mammal in need of such treatment with a compound of Claim 3.